

### Amendments to the Specification

Please amend the section beginning on page 59, line 21, as follows:

where R<sup>1</sup>, R<sup>4</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>64</sup>, Z and X are as defined in claim 15 and  
where X is O, or S, S(O) or S(O)<sub>2</sub>, NH or NR<sup>12</sup> where R<sup>12</sup> is hydrogen or C<sub>1-6</sub>alkyl;  
Z is C(O) or S(O)<sub>2</sub>.  
R<sup>1</sup> and R<sup>4</sup> are independently selected from halogeno, cyano, nitro, C<sub>1-3</sub>alkylsulphanyl,  
-N(OH)R<sup>13</sup>- (wherein R<sup>13</sup> is hydrogen, or C<sub>1-3</sub>alkyl), or R<sup>15</sup>X<sup>1</sup>- (wherein X<sup>1</sup> represents a direct  
bond, -O-, -CH<sub>2</sub>-, -OCO-, carbonyl, -S-, -SO-, -SO<sub>2</sub>-, -NR<sup>16</sup>CO-, -CONR<sup>16</sup>-,  
-NR<sup>17</sup>SO<sub>2</sub>- or -NR<sup>18</sup>- (wherein R<sup>16</sup>, R<sup>17</sup> and R<sup>18</sup> each independently represents hydrogen,  
C<sub>1-3</sub>alkyl or C<sub>1-3</sub>alkoxyC<sub>2-3</sub>alkyl), and R<sup>15</sup> is hydrogen, optionally substituted hydrocarbyl,  
optionally substituted heterocyclyl or optionally substituted alkoxy;  
R<sup>2</sup> and R<sup>3</sup> are groups R<sup>2</sup> and R<sup>3</sup> respectively, provided that at least one of said groups and  
preferably R<sup>3</sup> is a group of sub-formula X<sup>1</sup>-R<sup>15</sup> where X<sup>1</sup> is as defined above, and R<sup>15</sup> is a  
group R<sup>15</sup> as defined above in claim 4, provided that it is other than methyl, methyl;  
R<sup>6</sup> is hydrogen, optionally substituted hydrocarbyl or optionally substituted heterocyclyl;  
R<sup>7</sup> and R<sup>8</sup> are independently selected from hydrogen, halo, C<sub>1-4</sub>alkyl, C<sub>1-4</sub> alkoxy,  
C<sub>1-4</sub>alkoxymethyl, di(C<sub>1-4</sub>alkoxy)methyl, C<sub>1-4</sub>alkanoyl, trifluoromethyl, cyano, amino,  
C<sub>2-5</sub>alkenyl, C<sub>2-5</sub>alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic  
group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic  
group may be aromatic or non-aromatic and may be saturated (linked via a ring carbon or  
nitrogen atom) or unsaturated (linked via a ring carbon atom), and which phenyl, benzyl or  
heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected  
from hydroxy, halogeno, C<sub>1-3</sub>alkyl, C<sub>1-3</sub>alkoxy, C<sub>1-3</sub>alkanoyloxy, trifluoromethyl, cyano, amino,  
nitro, C<sub>2-4</sub>alkanoyl, C<sub>1-4</sub>alkanoylamino, C<sub>1-4</sub>alkoxycarbonyl, C<sub>1-4</sub>alkylsulphanyl,  
C<sub>1-4</sub>alkylsulphinyl, C<sub>1-4</sub>alkylsulphonyl, carbamoyl, N-C<sub>1-4</sub>alkylcarbamoyl,  
N,N-di(C<sub>1-4</sub>alkyl)carbamoyl, aminosulphonyl, N-C<sub>1-4</sub>alkylaminosulphonyl,  
N,N-di(C<sub>1-4</sub>alkyl)aminosulphonyl, C<sub>1-4</sub>alkylsulphonylamino, and a saturated heterocyclic  
group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl  
imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2  
substituents selected from oxo, hydroxy, halogeno, C<sub>1-3</sub>alkyl, C<sub>1-3</sub>alkoxy, C<sub>1-3</sub>alkanoyloxy,  
trifluoromethyl, cyano, amino, nitro and C<sub>1-4</sub>alkoxycarbonyl; and  
R<sup>64</sup> is optionally substituted hydrocarbyl or optionally substituted heterocyclyl.

Please amend the section beginning on page 61, line 13, as follows:

where X, Y, R<sup>1</sup>, R<sup>4</sup>, R<sup>7</sup>, R<sup>8</sup> are as defined in relation to compound (VIC), R<sup>65</sup> is as defined in relation to compound (VIC), and R<sup>68</sup> and R<sup>69</sup> are equivalent to R<sup>2</sup> and R<sup>3</sup> in relation to compound (VIC), except that at least one of R<sup>68</sup> or R<sup>69</sup> is a group of sub-formula X<sup>1</sup>R<sup>15</sup> where R<sup>15</sup> is as defined in relation to compound (VIC), provided that when said one of R<sup>68</sup> or R<sup>69</sup> is morpholinopropoxy, the other is not a group of sub-formula (18) as defined in claim 18 C<sub>2</sub>-alkenyl which may be unsubstituted or which may be substituted with one or more functional groups; and further provided that when said one of R<sup>68</sup> or R<sup>69</sup> is methoxyethoxy, the other is not methoxy.

Please amend the section beginning on page 64, line 18, as follows:

where X, R<sup>7</sup> and R<sup>8</sup> are as defined ~~in relation to the relevant compound according to any one of claims 19 to 26 for any one of formula (IIA), (IIB), (IIC), (IID), (VIA) or (VIB),~~ and R<sup>66</sup> is a group of formula NHZR<sup>64</sup> or Y(O)R<sup>65</sup> where Z, R<sup>64</sup>, Y and R<sup>65</sup> as are defined ~~in the relation to the said compound in any one of claims 19 to 26 for any one of formula (IIA), (IIB), (IIC), (IID), (VIA) or (VIB);~~ and thereafter if desired or necessary converting a group R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> or R<sup>4</sup> to a group R<sup>1</sup>, R<sup>2</sup> or R<sup>2</sup> or R<sup>68</sup>, R<sup>3</sup> or R<sup>3</sup> or R<sup>69</sup> and R<sup>4</sup> respectively or to a different such group.